Application No.: 10/622,320
Response dated March 6, 2006
Reply to Office Action of September 6, 2005

--Q₃ is a 5-6 membered aromatic carbocyclic or heterocyclic ring system, or an 8-10 membered bicyclic ring system comprising aromatic carbocyclic rings, aromatic heterocyclic rings or a combination of an aromatic carbocyclic ring and an aromatic heterocyclic ring. The rings of Q₃ are substituted with 1 to 4 substituents, each of which is independently selected from halo; C₁-C₃ alkyl optionally substituted with NR'₂, OR', CO₂R' or CONR'₂; O-(C₁-C₃)-alkyl optionally substituted with NR'₂, OR', CO₂R' or CONR'₂; NR'₂; OCF₃; CF₃; NO₂; CO₂R'; [[CONR']]CONHR'; SR'; S(O₂)N(R')₂; SCF₃; CN; N(R')C(O)R⁴; N(R')C(O)OR⁴; N(R')C(O)C(O)R⁴; N(R')S(O₂)R⁴; N(R')R⁴; N(R⁴)₂; OR⁴; OC(O)R⁴; OP(O)₃H₂; or [[N=C-N(R')₂]]N=CH-N(R')₂. --

Please replace the paragraph beginning at page 36, line 12, with the following rewritten paragraph:

--According to another preferred embodiment, Q₃ is a monocyclic carbocyclic ring, wherein each ortho substituent is independently selected from halo or methyl. According to another preferred embodiment, Q₃ contains 1 or 2 additional substituents independently selected from NR'₂, OR', CO₂R' CN, N(R')C(O)R⁴; N(R')C(O)OR⁴; N(R')C(O)C(O)R⁴; N(R')C(O)C(O)R⁴; N(R')C(O)C(O)R⁴; N(R')C(O)C(O)R⁴; N(R')C(O)C(O)R⁴; OP(O)₃H₂; or [[N=C-N(R')₂]]N=CH-N(R')₂.--Please replace the chemical structures of Table 4, depicted at pages 41-47, with

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--Table 4. Formula Ig Inhibitors.

the following redrawn chemical structures: